

- Hilberts space - abstract vector space with finite or infinite dimensionality

- Bra - Ket vectors: unit k vector usually

$|A\rangle + |B\rangle = |C\rangle$ $|A\rangle + |B\rangle = |B\rangle + |A\rangle$ $(|A\rangle + |B\rangle) + |C\rangle = |A\rangle + (|B\rangle + |C\rangle)$

$|A\rangle + (-|A\rangle) = 0$ $|zA\rangle = z|A\rangle$ $z(|A\rangle + |B\rangle) = z|A\rangle + z|B\rangle$

$(z+w)|A\rangle = z|A\rangle + w|A\rangle$

$z|A\rangle \Rightarrow \langle A|z^*$ $|A\rangle = \begin{pmatrix} \alpha_1 \\ \alpha_2 \\ \alpha_3 \\ \alpha_4 \end{pmatrix}$ $\langle A| = (\alpha_1^* \alpha_2^* \alpha_3^* \alpha_4^*)$

- inner product = $\langle B|A\rangle = \begin{pmatrix} \beta_1^* & \beta_2^* & \beta_3^* & \beta_4^* \end{pmatrix} \begin{pmatrix} \alpha_1 \\ \alpha_2 \\ \alpha_3 \\ \alpha_4 \end{pmatrix} = \alpha_1\beta_1^* + \alpha_2\beta_2^* + \alpha_3\beta_3^* + \alpha_4\beta_4^*$

$\langle B|A\rangle = \langle A|B\rangle^*$

$\langle C|(|A\rangle + |B\rangle) = \langle C|A\rangle + \langle C|B\rangle$

- normalized vector - of length 1 so gives inner product of 1

$\langle A|A\rangle = 1$ if A normalized

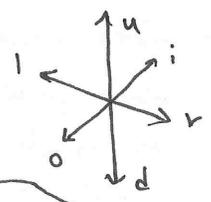
- orthogonal vectors - at right angles (perpendicular) to each other

$\langle B|A\rangle = 0$ if A, B orthogonal

- orthogonal normalized bases - orthonormal bases

- orthogonal + normalized used to orient in hilbert space

eg $\langle u|d\rangle = 0$
 $\langle l|r\rangle = 0$
 $\langle i|o\rangle = 0$



- linear operators - expressed by matrixes act on vectors

$M|A\rangle = |B\rangle$ $Mz|A\rangle = z|B\rangle$ $M(|A\rangle + |B\rangle) = M|A\rangle + M|B\rangle$

- eigen values and eigenvectors $|\lambda\rangle$ and λ $M|\lambda\rangle = \lambda|\lambda\rangle$

- Hermitian conjugation

$M|A\rangle = |B\rangle$ $\langle A|M^\dagger = \langle B|$ and $Mz|A\rangle = z|B\rangle$ $\langle A|z^*M^\dagger = \langle B|z^*$

M^\dagger - hermitian conjugate hermitian operator if $M = M^\dagger$

$M^\dagger = (M^T)^*$ $M = \begin{pmatrix} m_{11} & m_{12} & m_{13} \\ m_{21} & m_{22} & m_{23} \\ m_{31} & m_{32} & m_{33} \end{pmatrix}$

called transposing - reflect over diagonal so $m_{12} \Rightarrow m_{21}$

$M^T = \begin{pmatrix} m_{11} & m_{21} & m_{31} \\ m_{12} & m_{22} & m_{32} \\ m_{13} & m_{23} & m_{33} \end{pmatrix}$

find complex conjugate of each component so $m_{12} \Rightarrow m_{12}^*$

$M^\dagger = \begin{pmatrix} m_{11}^* & m_{12}^* & m_{13}^* \\ m_{21}^* & m_{22}^* & m_{23}^* \\ m_{31}^* & m_{32}^* & m_{33}^* \end{pmatrix}$

- probability of a state given state vector $|A\rangle$

$$|A\rangle = \alpha_u |u\rangle + \alpha_d |d\rangle$$

$$= \alpha_u \begin{pmatrix} 1 \\ 0 \end{pmatrix} + \alpha_d \begin{pmatrix} 0 \\ 1 \end{pmatrix}$$

$$= \begin{pmatrix} \alpha_u \\ \alpha_d \end{pmatrix} \quad \text{in terms of orthonormal vectors}$$

$$\alpha_u = \langle u|A\rangle$$

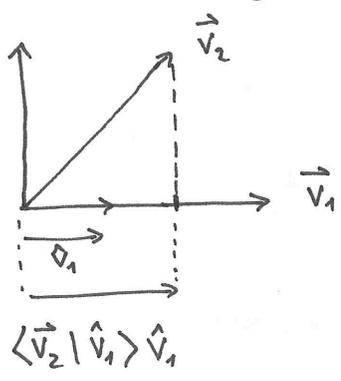
$$\alpha_d = \langle d|A\rangle$$

$$P_{(\text{spin up})} = P_u = \alpha_u^2 = \langle A|u\rangle \langle u|A\rangle$$

$$P_{(\text{spin down})} = P_d = \alpha_d^2 = \langle A|d\rangle \langle d|A\rangle$$

- Gram - Schmidt procedure

given \vec{v}_1 and \vec{v}_2 form orthonormal bases



$$\hat{v}_1 = \frac{\vec{v}_1}{|\vec{v}_1|}$$

$\langle \vec{v}_2 | \hat{v}_1 \rangle \hat{v}_1$ - component of \vec{v}_2 along \hat{v}_1
so "projected on \hat{v}_1 "

$\vec{v}_{2\perp}$ - component to \vec{v}_2 orthogonal to \hat{v}_1

$$\vec{v}_{2\perp} = \vec{v}_2 - \langle \vec{v}_2 | \hat{v}_1 \rangle \hat{v}_1 \quad \text{since } \vec{v}_2 = \vec{v}_{2\parallel} + \vec{v}_{2\perp}$$

parallel + normal

$$\hat{v}_2 = \frac{\vec{v}_{2\perp}}{|\vec{v}_{2\perp}|}$$

- observables are represented by Hermitian linear operators

the corresponding measured quantity is λ_i (eigenvalue)
and the state is $|\lambda_i\rangle$ so the eigenvector
 $|\lambda_i\rangle$ is the determined state by experiment

- state prepared in state $|A\rangle$ observing observable M

$$P(\lambda_i) = |\langle A | \lambda_i \rangle|^2$$

where $M|\lambda_i\rangle = \lambda_i|\lambda_i\rangle$

- spin represented by operator $\hat{S} = \begin{pmatrix} \hat{S}_x \\ \hat{S}_y \\ \hat{S}_z \end{pmatrix} = \begin{pmatrix} \sin\theta \cos\phi \\ \sin\theta \sin\phi \\ \cos\theta \end{pmatrix}$ in spherical coordinates

eg $\hat{S}_z |u\rangle$ and $|d\rangle$
 $\hat{S}_z = +1$ $\hat{S}_z = -1$

use to find θ and ϕ
then you can find other components such as \hat{S}_x, \hat{S}_y

$|u\rangle$ and $|d\rangle$ are eigenvectors of the spin operator

$$\hat{S}_z |u\rangle = (+1) |u\rangle \quad \hat{S}_z |d\rangle = (-1) |d\rangle$$

both are orthogonal $\langle u|d\rangle = 0$

eigen vectors:

$$|\lambda_1\rangle = \begin{pmatrix} \cos \frac{\theta}{2} \\ \sin \frac{\theta}{2} \end{pmatrix} \quad |\lambda_2\rangle = \begin{pmatrix} -\sin \frac{\theta}{2} \\ \cos \frac{\theta}{2} \end{pmatrix}$$

- pauli matrices for spin

$$\hat{\sigma}_x = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} \quad \hat{\sigma}_y = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix} \quad \hat{\sigma}_z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$$

- Spin polarization principle ^{single}

↳ any state of a spin is an eigenvector of some component of the spin

∴ given state $|A\rangle = \alpha_u|u\rangle + \alpha_d|d\rangle$

there exists a space direction $\hat{n} = \begin{pmatrix} n_x \\ n_y \\ n_z \end{pmatrix}$ for which $\vec{\sigma} \cdot \hat{n} |A\rangle = |A\rangle$

as a result $P_n = \langle \vec{\sigma} \cdot \hat{n} \rangle^2 = +1$ so $\vec{\sigma} \cdot \hat{n} = \lambda$ of $|A\rangle$
 $\lambda = +1$

↳ $\langle \sigma_x \rangle^2 + \langle \sigma_y \rangle^2 + \langle \sigma_z \rangle^2 = 1$ $\vec{\sigma} \cdot \hat{n} = \vec{\sigma}_n$ - component of spin $\vec{\sigma}$ along direction of \hat{n}

- Unitarity - condition that time evolution of quantum state according to Schrödinger's equation is represented by unitary operator

$|\psi(t)\rangle = U(t)|\psi(0)\rangle$ $U(t)$ - time development operator (time development operator)

- unitary operator $U^\dagger(t)U(t) = I$ I - unit operator

- states are distinguishable if they are orthogonal (there is an experiment to precisely tell them apart)

if $|\psi(0)\rangle, |\phi(0)\rangle$ are two orthogonal states therefore $\langle \psi(0) | \phi(0) \rangle = 0$

they will continue to be orthogonal for all t $\langle \psi(t) | \phi(t) \rangle = 0$

- characteristic equation for finding λ and $|\lambda\rangle$

$\det(M - \lambda I) = 0$ or $(M - \lambda I)|\lambda\rangle = 0$

$\det \begin{pmatrix} a & b \\ c & d \end{pmatrix} = ad - bc$

$I = \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}$ or $\begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{pmatrix}$

- time development also continuous so changes smoothly $\hbar \approx 1.05 \times 10^{-34} \text{ kg m}^2 \text{ s}^{-1}$

Hamiltonian - linear operator H

time independent Schrödinger equation: $\hbar \frac{\partial |\psi\rangle}{\partial t} = -iH|\psi\rangle$

- average or expected value $\langle L \rangle = \langle A | L | A \rangle$ where L - linear operator
 $= \sum_i (\alpha_i^* \alpha_i) \lambda_i$ $|A\rangle$ - state vector

these factors can be ignored as they cancel out expectation value at time t
 $\langle L \rangle = \langle A | e^{-iHt} L e^{iHt} | A \rangle$
 $= \langle A | L e^0 | A \rangle$
 $= \langle A | L | A \rangle$
 $= \langle \psi(t) | L | \psi(t) \rangle$

- commutation: - commutator of L with M:

$$[L, M] = LM - ML \quad [L, M] = -[M, L]$$

$$\frac{d\langle E \rangle}{dt} = -\frac{i}{\hbar} \langle [L, H] \rangle \quad \frac{dL}{dt} = -\frac{i}{\hbar} [L, H]$$

symmetrical to classical mechanics
where $\frac{dF}{dt} = \{F, H\}$

omitting the angle brackets $\langle \rangle$

- converting quantum commutators to poisson brackets: $[F, G] = i\hbar \{F, G\}$

now we can do

in this case $[L, H] = i\hbar \{L, H\}$

$$\frac{dL}{dt} = -\frac{i}{\hbar} (i\hbar \{L, H\}) = \{L, H\}$$

where L is an observable

- conservation

if quantity or operator commutes with hamiltonian H it is conserved $[Q, H] = 0$ this works for all powers of Q so Q^n

therefore ~~if quantity~~ all functions of Q^n conserved

↳ if $[Q^n, H] = 0$ then quantity tied to observable Q then it is conserved

eg H - represents energy $[H, H] = 0 \therefore$ Energy conserved

- spin in the magnetic field

energy proportional dot product of spin and magnetic field

$$H \propto \vec{\sigma} \cdot \vec{B} = \sigma_x B_x + \sigma_y B_y + \sigma_z B_z$$

field lies along z axis so $B_x = 0 \quad B_y = 0 \therefore H \propto \sigma_z B_z$

$$H = \frac{\hbar \omega}{2} \sigma_z \text{ where } \omega \text{ contains all constants for simplicity}$$

$$\langle \dot{\sigma}_x \rangle = -\frac{i}{\hbar} \langle [\sigma_x, H] \rangle = -\frac{i\omega}{2} \langle [\sigma_x, \sigma_z] \rangle = -\omega \langle \sigma_y \rangle$$

$$\langle \dot{\sigma}_y \rangle = -\frac{i}{\hbar} \langle [\sigma_y, H] \rangle = -\frac{i\omega}{2} \langle [\sigma_y, \sigma_z] \rangle = \omega \langle \sigma_x \rangle$$

$$\langle \dot{\sigma}_z \rangle = -\frac{i}{\hbar} \langle [\sigma_z, H] \rangle = -\frac{i\omega}{2} \langle [\sigma_z, \sigma_z] \rangle = 0 \text{ since they commute}$$

use pauli matrices for spin to compute commutators

$$\text{so } [\sigma_x, \sigma_z] = 2i\sigma_y \quad [\sigma_y, \sigma_z] = 2i\sigma_x \quad [\sigma_z, \sigma_z] = 2i\sigma_y$$

again $x \rightarrow y \rightarrow z$

- solving Schrödinger equation

$$i\hbar \frac{\partial \Psi(x)}{\partial t} = -\frac{\hbar^2}{2m} \frac{\partial^2 \Psi(x)}{\partial x^2} + U(x) \Psi(x)$$

↳ procedure:

1. find and express Hamiltonian $H =$

2. prepare an initial state $|\Psi(0)\rangle$

3. solve time independent equation to get λ_j and $|\lambda_j\rangle$
 $H|E_j\rangle = E_j|E_j\rangle$ or E_j and $|E_j\rangle$

4. use $|\Psi(0)\rangle$ and $E_j, |E_j\rangle$ to calculate initial coefficients $\alpha_j(0)$
 $\alpha_j(0) = \langle E_j | \Psi(0) \rangle$

5. rewrite $|\Psi(0)\rangle$ in terms of eigenvectors $|E_j\rangle$ and initial coefficients $\alpha_j(0)$

$$|\Psi(0)\rangle = \sum_j \alpha_j(0) |E_j\rangle \quad \text{check if still makes sense}$$

6. replace initial with time dependant $0 \Rightarrow t$

$$|\Psi(t)\rangle = \sum_j \alpha_j(t) |E_j\rangle \quad \left. \begin{array}{l} \\ \end{array} \right\} \text{eigenvalue of } |E_j\rangle$$

7. using equation below express $\alpha_j(t)$ in terms of α_j, E_j and t

~~$$\alpha_j(0) = \langle E_j | \Psi(0) \rangle$$~~ so $|\Psi(t)\rangle = \sum_j \alpha_j(0) e^{-\frac{i}{\hbar} E_j t}$

$$\alpha_j(t) = \alpha_j(0) e^{-\frac{i}{\hbar} E_j t}$$

8. values can be predicted using eigenvalues and eigenvectors

suppose observing observable L with eigenvalue λ and vector $|\lambda\rangle$

$$P_\lambda(t) = |\langle \lambda | \Psi(t) \rangle|^2 = \langle \lambda | \Psi(t) \rangle \langle \Psi(t) | \lambda \rangle$$

- solving time dependant Schrödinger equation short:

1. find or get Hamiltonian

2. prepare initial state $|\Psi(0)\rangle$

3. solve $H|E_j\rangle = E_j|E_j\rangle$ so find eigenvalue of state vectors

4. find initial coefficients $\alpha_j(0) = \langle E_j | \Psi(0) \rangle$

5. substitute into $|\Psi(0)\rangle = \sum_j \alpha_j(0) |E_j\rangle$ α_j is coefficient $|E_j\rangle$ is state vector

6. solve $\alpha_j(t) = \alpha_j(0) e^{-\frac{i}{\hbar} E_j t}$

7. substitute into $|\Psi(t)\rangle = \sum_j \alpha_j(t) |E_j\rangle$

8. simplify if necessary. check using $t=0$ if $|\Psi(0)\rangle$ is obtained.

- note - a measurement leaves the system in an eigenstate (state of a single eigenvector) corresponding to the last value (eigenvalue) measured

- eg if λ_i measured state is now $|\lambda_i\rangle$

- state of multiple measurables

↳ consider 2 independent spins. each has an observable, eigenvalue, eigenvector corresponding to hermitian operator L and M

after a measurement system is in a state which is both an eigenvector of L and M called simultaneous eigenvector of operators

state vector of each spin: $L|\lambda\rangle = \lambda|\lambda\rangle$ $M|\mu\rangle = \mu|\mu\rangle$

bases of state vectors which are simultaneous eigenvectors of both L, M

$LM|\lambda, \mu\rangle = \lambda\mu|\lambda, \mu\rangle$ in order for this to be possible L, M must commute

since order of multiplication of eigenvalues doesn't matter

∴ $LM|\lambda, \mu\rangle = ML|\lambda, \mu\rangle$ so $[L, M]|\lambda, \mu\rangle = [L, M]|\lambda, \mu\rangle = 0$ since $[L, M] = 0$

- the condition for 2 observables to be simultaneously measured is for their hermitian operators to commute

- if 2 hermitian operators commute there exists a complete basis of simultaneous eigenvectors (and eigenvalues) of the observables

- for higher n of observables they must all commute with each other they are called complete set of commuting observables

- wave function

suppose bases for some quantum system we have: - orthonormal basis vectors $|a, b, c, \dots\rangle$

since $|a, b, c, \dots\rangle$ orthonormal

$|\Psi\rangle$ can be expressed in terms of $|a, b, c, \dots\rangle$

$|\Psi\rangle = \sum_{a,b,c,\dots} \psi(a,b,c,\dots) |a,b,c,\dots\rangle$

- eigenvalues a, b, c, \dots

- complete set of commuting observables A, B, C, \dots

- arbitrary state vector $|\Psi\rangle$

$\psi(a, b, c, \dots)$ are coefficients for expansion

also each coefficient = inner product of $|\Psi\rangle$ with basis vector

so $\psi(a) = \langle a | \Psi \rangle$ for all: $\psi(a, b, c, \dots) = \langle a, b, c, \dots | \Psi \rangle$

$\psi(a, b, c, \dots)$ - is the wave function

$P(a) = \psi^*(a) \psi(a)$ for all $P(a, b, c, \dots) = \psi^*(a, b, c, \dots) \psi(a, b, c, \dots)$

wave function in δz $\psi(u) = \langle u | \Psi \rangle$ $\psi(d) = \langle d | \Psi \rangle$

P must always sum to 1

$\sum_{a,b,c,\dots} \psi^*(a,b,c,\dots) \psi(a,b,c,\dots) = 1$

$\sum_i \psi^*(i) \psi(i) = 1$

- Any 2x2 Hermitian matrix representing operator can be written as sum of 4 terms with real coefficients

$$L = a\delta_x + b\delta_y + c\delta_z + d\delta I \quad a, b, c, d \in \mathbb{R} \quad \delta_x, \delta_y, \delta_z - \text{Pauli matrices}$$

- Spin components can not be measured simultaneously since don't commute

$$[\delta_x, \delta_y] = 2i\delta_z \quad [\delta_y, \delta_z] = 2i\delta_x \quad [\delta_z, \delta_x] = 2i\delta_y$$



$$\therefore [\delta_x, \delta_y] \neq 0$$

- Uncertainty principle

↳ if hermitian operators of 2 observables do not commute there must be uncertainty in measurement of one/other/both

given state $|\psi\rangle$ and observable A with eigenvalue a

$$P(a) \text{ is probability of observing } a \quad \langle \psi | A | \psi \rangle = \sum_a a P(a)$$

since expected val = $\sum \text{val} \times \text{probability of value}$

- uncertainty in A is its standard deviation

$$\text{begin by defining } \bar{A} \quad \bar{A} = A - \langle \psi | A | \psi \rangle I \text{ so } \bar{A} = A - \text{expectation value} \times I$$
$$= A - \langle A \rangle I$$

we use I as unit operator to subtract $\langle A \rangle$ which is a single \mathbb{R} value

$$\text{this also shifts eigenvalues of } A \quad \bar{a} = a - \langle \psi | A | \psi \rangle I$$
$$= a - \langle A \rangle I$$

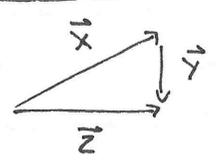
$$\text{if uncertainty} = \text{standard deviation} = \Delta A \quad \text{variance} = (\Delta A)^2$$

$$(\Delta A)^2 = \sum_a \bar{a}^2 P(a) = \sum_a (a - \langle A \rangle I)^2 P(a)$$

$$\text{which can be written as } (\Delta A)^2 = \langle \psi | \bar{A}^2 | \psi \rangle \quad \text{if } \langle A \rangle = 0 \quad (\Delta A)^2 = \langle \psi | A^2 | \psi \rangle$$

- Cauchy-Schwarz Inequality

in real vector space $|x| + |y| \geq |x+y|$
which gives $|x||y| \geq |x \cdot y|$



$|x+y|$ is shortest $|z|$ can be

↳ length of vector = its square root of its inner product with itself

$$|\vec{M}| = \sqrt{\langle M | M \rangle}$$

$$\text{Cauchy-Schwarz inequality: } |\vec{x}| |\vec{y}| \geq |\vec{x} \cdot \vec{y}|$$

↳ the form used in uncertainty principle:

$$2|\vec{x}| |\vec{y}| \geq |\langle x | y \rangle + \langle y | x \rangle|$$

- General Uncertainty principle

let $|\psi\rangle$ be any state A, B are any 2 observables

define $|x\rangle = A|\psi\rangle$ $|y\rangle = iB|\psi\rangle$ (don't forget i!)

substitute $|x\rangle$ and $|y\rangle$ into Cauchy Schwarz inequality

$$2|\langle x|y\rangle| \geq |\langle x|x\rangle| |\langle y|y\rangle|$$

$$2\sqrt{\langle A^2\rangle} \sqrt{\langle B^2\rangle} \geq |\langle \psi|AB|\psi\rangle - \langle \psi|BA|\psi\rangle|$$

$$2\sqrt{\langle A^2\rangle} \sqrt{\langle B^2\rangle} \geq |\langle \psi|[A,B]|\psi\rangle| \quad \text{using notation of commutators}$$

$$\Delta \bar{A} \Delta \bar{B} \geq \frac{1}{2} |\langle \psi|[A, B]|\psi\rangle|$$

$$\Delta A \Delta B \geq \frac{1}{2} |\langle \psi|[A, B]|\psi\rangle|$$

product of uncertainties can not be smaller than $\frac{1}{2}$ magnitude of expectation value of commutator of observables

Heisenberg's uncertainty principle
 ↳ product of uncertainty of position (q) and momentum (p) is greater or equal to $\frac{1}{2}$ Planck's constant ($\frac{1}{2} \hbar$)

- states of a combined system

2 independent spins Alice and Bob can have states $|u\rangle$ and $|d\rangle$ each

they exist in state space S_A and S_B

we define the state space of both Alice and Bob S_{AB}

$$S_{AB} = S_A \otimes S_B \quad \text{where } \otimes \text{ is the tensor product}$$

the orthonormal basis of S_{AB} are $|uu\rangle, |ud\rangle, |du\rangle, |dd\rangle$

- statistical correlation - shows how correlated observations of Alice and Bob are

if Alice and Bob toss a coin $H=+1$ $T=-1$ averages $\langle \delta_A \rangle = 0$

if system is correlated so for every H Alice gets a T $\langle \delta_B \rangle = 0$

the product $\delta_A \delta_B = -1$ allways so $\langle \delta_A \delta_B \rangle = -1$

this would be perfect (100%) correlation

$$\text{statistical correlation} = \langle \delta_A \delta_B \rangle - \langle \delta_A \rangle \langle \delta_B \rangle$$

if 2 variables completely uncorrelated $P(a,b) = P_A(a) P_B(b)$ so factorise

- Product states

$$|\text{product state}\rangle = (\alpha_u |u\rangle + \alpha_d |d\rangle) \otimes (\beta_u |u\rangle + \beta_d |d\rangle)$$

$$= \alpha_u \beta_u |uu\rangle + \alpha_u \beta_d |ud\rangle + \alpha_d \beta_u |du\rangle + \alpha_d \beta_d |dd\rangle$$

if state vectors for Alice ($|u\rangle, |d\rangle$) and Bob ($|u\rangle, |d\rangle$)

are normalized the product state is automatically normalized as well

it took 2 \mathbb{R} parameters for a single ^{spin} state system now it is 4 \mathbb{R} parameters

$$(\alpha_u, \alpha_d, \beta_u, \beta_d)$$

- entanglement is not "all or nothing" proposition there are levels of entanglement

- singlet and triplet states are most entangled states possible and can not be written in product ~~form~~ state

$$|sing\rangle = \frac{1}{\sqrt{2}} (|ud\rangle - |du\rangle) \quad |T_1\rangle = \frac{1}{\sqrt{2}} (|ud\rangle + |du\rangle)$$

$$|T_2\rangle = \frac{1}{\sqrt{2}} (|uu\rangle + |dd\rangle) \quad |T_3\rangle = \frac{1}{\sqrt{2}} (|duu\rangle - |dd\rangle)$$

- an entangled state gives complete description of combined system no more can be known

- in maximally entangled state nothing is known about the individual systems even if you know entangled state

- Tensor product operators

$$\sigma_z |du\rangle = -|du\rangle$$

$$(\sigma_z \otimes I) (|d\rangle \otimes |u\rangle) = (\sigma_z |d\rangle \otimes I |u\rangle)$$

$$= (-|d\rangle \otimes |u\rangle)$$

I is unit operator so σ_z only affects ~~both~~ alices (first) half likewise T_z would affect bobs half only

- spin-polarization principle - not all expectation values can be 0

$$\langle \sigma_x \rangle^2 + \langle \sigma_y \rangle^2 + \langle \sigma_z \rangle^2 = 1$$

applies to all entangled states except singlet state $|sing\rangle$

sing goes to 0 for $|sing\rangle \quad \langle \sigma_x \rangle = \langle \sigma_y \rangle = \langle \sigma_z \rangle = 0$

$$\vec{\sigma} \cdot \vec{T} = \sigma_x T_x + \sigma_y T_y + \sigma_z T_z$$

$$= (\sigma_x \otimes I)(I \otimes T_x) + (\sigma_y \otimes I)(I \otimes T_y) + (\sigma_z \otimes I)(I \otimes T_z)$$

- for tensor products

$$\alpha = \begin{pmatrix} a_1 & a_2 \\ a_3 & a_4 \end{pmatrix} \quad \beta = \begin{pmatrix} \beta_1 & \beta_2 \\ \beta_3 & \beta_4 \end{pmatrix}$$

$$\alpha \otimes \beta = \begin{pmatrix} a_1 \begin{pmatrix} \beta_1 & \beta_2 \\ \beta_3 & \beta_4 \end{pmatrix} & a_2 \begin{pmatrix} \beta_1 & \beta_2 \\ \beta_3 & \beta_4 \end{pmatrix} \\ a_3 \begin{pmatrix} \beta_1 & \beta_2 \\ \beta_3 & \beta_4 \end{pmatrix} & a_4 \begin{pmatrix} \beta_1 & \beta_2 \\ \beta_3 & \beta_4 \end{pmatrix} \end{pmatrix}$$

if α has N_α dimensions and β has N_β dimensions

$\alpha \otimes \beta$ has $N_\alpha N_\beta$

eg $\alpha: 3 \times 4 \quad \beta: 2 \times 5$

$$\alpha \otimes \beta: 12 \times 10 = 120$$

- can be applied for all ev soly schrodinger the detent

- given matrix M and basis vectors j, k we can calculate the matrix entries of M_{jk} with respect to the basis

$$M_{jk} = \langle j | M | k \rangle \quad \text{eg) } M = \begin{pmatrix} 1 & 2 \\ 3 & 4 \end{pmatrix} \quad M_{jk} = \langle j | M | k \rangle = (1 \ 0) \begin{pmatrix} 1 & 2 \\ 3 & 4 \end{pmatrix} \begin{pmatrix} 1 \\ 0 \end{pmatrix} \\ = (1 \ 2) \begin{pmatrix} 1 \\ 0 \end{pmatrix} = 1 \quad \left(\begin{matrix} 1 \\ 0 \end{matrix} \right)$$

- Tensor product matrices:

$$A \otimes B = \begin{pmatrix} A_{11}B & A_{12}B \\ A_{21}B & A_{22}B \end{pmatrix} = \begin{pmatrix} A_{11}B_{11} & A_{11}B_{12} & A_{12}B_{11} & A_{12}B_{12} \\ A_{11}B_{21} & A_{11}B_{22} & A_{12}B_{21} & A_{12}B_{22} \\ A_{21}B_{11} & A_{21}B_{12} & A_{22}B_{11} & A_{22}B_{12} \\ A_{21}B_{21} & A_{21}B_{22} & A_{22}B_{21} & A_{22}B_{22} \end{pmatrix} \quad \text{eg) } |uu\rangle = \begin{pmatrix} 1 \\ 0 \end{pmatrix} \otimes \begin{pmatrix} 1 \\ 0 \end{pmatrix} \\ = \begin{pmatrix} 1 \\ 0 \\ 0 \\ 0 \end{pmatrix}$$

- outer product gives linear operator

outer product: $|\psi\rangle\langle\phi|$ if it acts on $|A\rangle$: $|\psi\rangle\langle\phi|A\rangle \equiv |\psi\rangle\langle\phi|A\rangle$
 if it acts on $\langle B|$: $\langle B| |\psi\rangle\langle\phi| \equiv \langle B|\psi\rangle\langle\phi|$

- outer product of same bra and ket = projection operator

$|\psi\rangle\langle\psi|$ projects vector onto direction of $|\psi\rangle$

- projection operators: - are Hermitian $(|\psi\rangle\langle\psi|)^\dagger = |\psi\rangle\langle\psi|$

- any vector orthogonal to $|\psi\rangle$ is an eigenvector with eigenvalue of 0
- vector $|\psi\rangle$ is an eigenvector with eigenvalue of 1
- square of projection operator is itself $(|\psi\rangle\langle\psi|)^2 = |\psi\rangle\langle\psi|$
- ~~trace of an operator is 1~~

- Trace of the projection operator is 1 $\text{Tr } |\psi\rangle\langle\psi| = 1$

$\sum_i |i\rangle\langle i| = I \longrightarrow$ - sum of projection operators of basis system gives Identity operator I

- Trace is defined as the sum of matrices diagonal entries

$$\text{eg) } \text{Tr} \begin{pmatrix} 6 & 5 & 3 \\ 4 & 2 & 0 \\ 1 & 0 & 5 \end{pmatrix} = 6 + 2 + 5 = 13$$

- Trace of a hermitian operator is the sum of its eigenvalues

$$\text{Tr } M = \sum_i \lambda_i$$

- expectation value of any observable L in state $|\psi\rangle$ can be calculated as:
 $\langle\psi|L|\psi\rangle = \text{Tr } |\psi\rangle\langle\psi|L$

- Density matrices

if alice prepares state $|\psi\rangle$ expectation value = $\text{Tr } |\psi\rangle\langle\psi|L$

if alice prepares state $|\phi\rangle$ expectation value = $\text{Tr } |\phi\rangle\langle\phi|L$

if 50% chce $|\psi\rangle$ and 50% $|\phi\rangle$ expectation value = $\frac{1}{2} \text{Tr } |\psi\rangle\langle\psi|L + \frac{1}{2} \text{Tr } |\phi\rangle\langle\phi|L$
 and density matrix would be $\rho = \frac{1}{2} |\psi\rangle\langle\psi| + \frac{1}{2} |\phi\rangle\langle\phi|$

so expectation value $\langle L \rangle = \text{Tr } \rho L$

- density matrix $\rho = P_1 |\phi_1\rangle\langle\phi_1| + P_2 |\phi_2\rangle\langle\phi_2| + P_3 |\phi_3\rangle\langle\phi_3| + \dots$

$$\rho = \sum_i (P_i |\phi_i\rangle\langle\phi_i|)$$

density matrix is an operator until basis are chosen

eg) $\rho_{aa'} = \langle a | \rho | a' \rangle$ $\langle L \rangle = \sum_{aa'} L_{a',a} \rho_{a,a'}$

- density matrices and entanglement

to calculate Alices density matrix ρ we might need the full wave function $\Psi(a,b)$ including bobs half. once we know ρ we can forget where we obtained it from and use it to calculate anything about Alices observations

eg to calculate $P(a)$ so alices state to be a start $P(a,b) = \Psi^*(a,b) \Psi(a,b)$ by rules of probability summing over b gives $P(a)$

$$P(a) = \sum_b P(a,b) = \sum_b \Psi^*(a,b) \Psi(a,b) \rightarrow \text{this is the diagonal entry in the density matrix} \\ = \rho_{aa}$$

- properties of density matrices

↳ Hermitian $\rho_{aa'} = \rho_{a'a}^*$ $(\rho_{aa'}^T = \rho_{a'a})$ transpose just exchange entries in subscript

↳ Trace of density matrix is 1 $\text{Tr}(\rho) = 1$ (sum of probabilities must = 1)

↳ eigenvalues of ρ are all +ve and lie between 0 and 1
it follows if one $\lambda = 1$ others must be 0
(since sum of probability = 1 and one probability is 1 others = 0)

↳ for a pure state $\rho^2 = \rho$ $\text{Tr}(\rho^2) = 1$ (pure - unentangled)

↳ for a mixed state $\rho^2 \neq \rho$ $\text{Tr}(\rho^2) < 1$ (mixed - entangled)

- Trace of product of 2 matrices is commutative

$$\text{Tr}(AB) = \text{Tr}(BA) \quad AB \neq BA$$

- Calculating alices ρ (u and d are 2 states for alice)

$$\rho_{a'a} = \sum_b \Psi^*(a,b) \Psi(a',b) \quad \text{consider state vector } |\Psi\rangle = \frac{1}{\sqrt{2}} (|ud\rangle + |du\rangle)$$

recall basis vector coefficients $\Psi(u,u) = 0$ $\Psi(u,d) = \frac{1}{\sqrt{2}}$
 $\Psi(d,d) = 0$ $\Psi(d,u) = \frac{1}{\sqrt{2}}$

set $b = b'$ in situation and find entries of ρ

$$\rho_{uu} = \Psi^*(u,u) \Psi(u,u) + \Psi^*(u,d) \Psi(u,d) = \frac{1}{2}$$

$$\rho_{ud} = \Psi^*(d,u) \Psi(u,u) + \Psi^*(d,d) \Psi(u,d) = 0$$

$$\rho_{du} = \Psi^*(u,u) \Psi(d,u) + \Psi^*(u,d) \Psi(d,d) = 0$$

$$\rho_{dd} = \Psi^*(d,u) \Psi(d,u) + \Psi^*(d,d) \Psi(d,d) = \frac{1}{2}$$

$$\rho = \begin{pmatrix} \rho_{uu} & \rho_{ud} \\ \rho_{du} & \rho_{dd} \end{pmatrix} = \begin{pmatrix} \frac{1}{2} & 0 \\ 0 & \frac{1}{2} \end{pmatrix} = \frac{1}{2} I$$

check $\text{Tr}(\rho) = 1$

- Test for entanglement - if entangled the wave function $\Psi(a,b)$ and probability function $P(a,b)$ will not factorise \therefore correlated

or assume A is alices observable B is bobs observable with corresponding expectation values $\langle A \rangle, \langle B \rangle$

correlation between them $C(A,B) = \langle AB \rangle - \langle A \rangle \langle B \rangle$

if $C(A,B) \neq 0$ state is entangled if $C(A,B) = 0$ so $\langle AB \rangle = \langle A \rangle \langle B \rangle$ than pure state
 $-1 \leq C(A,B) \leq +1$

- for product state (not entangled) the density matrix has only 1 non-zero eigen value equal to exactly 1 $\therefore \sum_a P a' a \alpha_a = \lambda \alpha_a' \quad \lambda = 1$
the eigen vector of this eigen value of 1 will be the wave function Ψ

- for singlet state or any maximally entangled the density matrix is ~~equal~~ to proportional to unit matrix since all eigen values sum up to I (1) eigen value is $\frac{1}{N_A}$
 N_A - alices state of states
- number of alices states
 $P a' a = \lambda \delta_{a' a}$
 $= \frac{1}{N_A} \delta_{a' a}$ Kronecker delta so probability of all states is equal

- Measurement

take apparatus A with states $|b\rangle, |+1\rangle, |-1\rangle$ spin states $|u\rangle, |d\rangle$
b is blank
+1 is up
-1 is down

both have states since both are quantum states using tensor product create space of states
if spin initially in $|u\rangle$ and A in $|b\rangle$ they interact to give A in $|+1\rangle$ so $|u, b\rangle \rightarrow |u, +1\rangle$
 $|u, b\rangle$
 $|u, +1\rangle$
 $|u, -1\rangle$
 $|d, b\rangle$
 $|d, +1\rangle$
 $|d, -1\rangle$

generally initial state of spin $\alpha_u |u\rangle + \alpha_d |d\rangle$ and of system with A $\alpha_u |u, b\rangle + \alpha_d |d, b\rangle$

which gives final state $\alpha_u |u, +1\rangle + \alpha_d |d, -1\rangle$
so $\alpha_u |u, b\rangle + \alpha_d |d, b\rangle \rightarrow \alpha_u |u, +1\rangle + \alpha_d |d, -1\rangle$

final state is entangled so A entangled with measured system (must be entangled because from state of A we can tell spin)

expectation value of +1 is still $\alpha_u^* \alpha_u$ until alice looks at it but for bob she did not collapse the wave function as he didnt look at system yet she just became entangled with the system if bob looks at her he is also entangled with system but there is still charlie...

- entanglement + locality
 - ↳ John Bell proved QM is non local
 - ↳ locality - impossible to send signal (information) faster than light
 - ↳ even for maximally entangled state Bob can do nothing to immediately affect alices wave function or density matrix
 - so one could argue QM is local this is ensured by the operator acting being unitary $U^\dagger = U$

- Bells theorem - any classical simulation ~~which~~ of QM which tries to spatially separate alices and bobs apparatuses and systems must have an instantaneous "cable" connecting the two with central memory to update the state vectors (cable only to simulate entanglement)
 - in real world no cable needed ~~since proven to be~~ but in simulation on boolean computer yes
 - information sent faster than speed of light but not accessed faster so cable even in real life but can not be discovered we know it exists since we need to include it in our simulations.

Summary:

- Product state - no entanglement \therefore excessive locality, seems as classical system
 - each sub system is fully characterized. no correlation between alices and bobs system
 - state vector: $\alpha_u \beta_u |uu\rangle + \alpha_u \beta_d |ud\rangle + \alpha_d \beta_u |du\rangle + \alpha_d \beta_d |dd\rangle$
 - normalization: $\alpha_u^* \alpha_u + \alpha_d^* \alpha_d = 1 \quad \beta_u^* \beta_u + \beta_d^* \beta_d = 1$
 - density matrix has 1 nonzero λ and is = to 1 the eigen vector is the alices wave function same for bob
 - wave function factorised $\Psi(a) \phi(b)$
 - expectation values $\langle \sigma_x \rangle^2 + \langle \sigma_y \rangle^2 + \langle \sigma_z \rangle^2 = 1 \quad \langle \tau_x \rangle^2 + \langle \tau_y \rangle^2 + \langle \tau_z \rangle^2 = 1$
 - correlation $\langle \sigma_z \tau_z \rangle - \langle \sigma_z \rangle \langle \tau_z \rangle = 0$
- Singlet state - maximum entangled \therefore non locality with all quantum weirdness
 - composite system fully characterised but no information about alices or bobs sub systems
 - state vector: $\frac{1}{\sqrt{2}} (|ud\rangle - |du\rangle)$
 - normalization: $\Psi_{uu}^* \Psi_{uu} + \Psi_{ud}^* \Psi_{ud} + \Psi_{du}^* \Psi_{du} + \Psi_{dd}^* \Psi_{dd} = 1$
 - density matrix has full composite system $\rho^2 = \rho \quad \text{Tr}(\rho^2) = 1$
alices density matrix is proportional to \mathbb{I} with equal eigen values which all add up to 1 so all measured outcomes are equally likely for subsystem of alice $\rho^2 \neq \rho \quad \text{Tr}(\rho^2) \neq 1$
 - wave function not factorised - $\Psi(a, b)$
 - expectation values $\langle \sigma_x \rangle = \langle \sigma_y \rangle = \langle \sigma_z \rangle = 0 \quad \langle \tau_x \rangle = \langle \tau_y \rangle = \langle \tau_z \rangle = 0$

correlation: $\langle \sigma_z \tau_z \rangle - \langle \sigma_z \rangle \langle \tau_z \rangle = -1$
 $\langle \tau_z \sigma_z \rangle, \langle \tau_x \sigma_x \rangle, \langle \tau_y \sigma_y \rangle = -1$

- Almost singlet state - partial entanglement - difficult to tell up-down
 - some information about composite system some about individual systems. In both cases incomplete
 - state vector: $\sqrt{0.6}|ud\rangle - \sqrt{0.4}|du\rangle$
 - normalization: $\Psi_{uu}^* \Psi_{uu} + \Psi_{ud}^* \Psi_{ud} + \Psi_{du}^* \Psi_{du} + \Psi_{dd}^* \Psi_{dd} = 1$
- density matrix of
 - ↳ full composite system $\rho^2 = \rho$ $\text{Tr}(\rho^2) = 1$
 - ↳ Alice's subsystem $\rho^2 \neq \rho$ $\text{Tr}(\rho^2) < 1$
- wave function not factorised $\Psi(a, b)$
- expectation values $\langle \delta_z \rangle = 0.2$ $\langle \delta_x \rangle = \langle \delta_y \rangle = 0$
 $\langle T_z \rangle = -0.2$ $\langle T_x \rangle = \langle T_y \rangle = 0$
 $\langle T_z \delta_z \rangle = -1$ $\langle T_x \delta_x \rangle = -2\sqrt{0.24}$
- correlation $\langle \delta_z T_z \rangle - \langle \delta_z \rangle \langle T_z \rangle = -0.56$
 otherwise $-1 < C(a, b) < 1$ $C(a, b) \neq 0$ for partially entangled

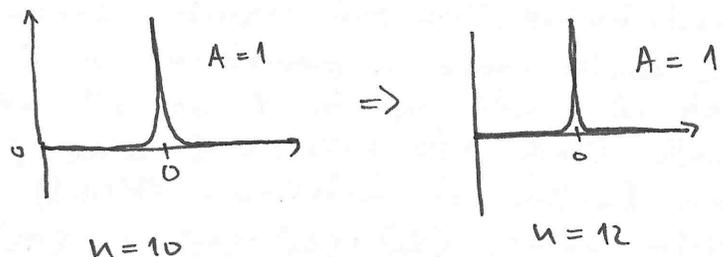
- continuous wave functions
 - ↳ sums \Rightarrow integrals
 - ↳ probability \Rightarrow probability density
 - ↳ Kronecker delta \Rightarrow Dirac delta

- probability density with discrete wave functions $P(x) = \Psi^*(x)\Psi(x)$
 but with continuous it is near impossible for it to be exactly x so we find P of a range of values $P(a, b) = P(a < x < b)$

- so we define a normalized vector by $\int_{-\infty}^{\infty} \Psi^*(x)\Psi(x) dx = 1$

- Dirac delta function
 - ↳ it is zero for all values $x \neq x'$ but if $x = x'$ it has infinite magnitude. this occurs at an infinitesimal range ϵ and the amplitude is $\frac{1}{\epsilon}$ so the area is 1

$\frac{n}{\sqrt{\pi}} e^{-(nx)^2}$
 as $n \rightarrow \infty$



as $n \uparrow$ width decreases

- linear operator - "machine" which acts on a function to give another function

- ↳ linear ∴ - if it acts on ~~the function~~ sum of 2 functions it gives sum of individual results
- if it acts on multiple of function it gives multiple of results

e.g operator X: $X\psi(x) = x\psi(x)$

D: $D\psi(x) = \frac{d\psi(x)}{dx}$

- for hermitian operators $L^\dagger = L$ it also holds that then between bra and kets are complex conjugates

$$\langle \Psi | L | \Phi \rangle = (\langle \Phi | L | \Psi \rangle)^*$$

e.g.) $\langle \Psi | X | \Phi \rangle = \int \psi^*(x) x \phi(x) dx$ $\langle \Phi | X | \Psi \rangle = \int x \psi(x) \phi^*(x) dx$

D is not hermitian but anti-hermitian for which

$$D^\dagger = -D$$

we correct this by a constant $-i\hbar$ so we get $-i\hbar D$

- eigen values and eigenvectors of position

$X|\Psi\rangle = x_0|\Psi\rangle$ so $x\psi(x) = x_0\psi(x)$ we get $\langle x|\Psi\rangle = \psi(x)$

so wave function is projection of state vector $|\Psi\rangle$ onto the eigenvector $|x\rangle$

- eigen values and eigenvectors of momentum

$P = -i\hbar D$ $P\psi(x) = -i\hbar \frac{d\psi(x)}{dx}$ eigen equation: $P|\Psi\rangle = p|\Psi\rangle$

$$\frac{d\psi(x)}{dx} = \frac{ip}{\hbar} \psi(x)$$

general solution $\psi_p(x) = A e^{\frac{ipx}{\hbar}}$

↳ specific eigenvector of P with eigenvalue p

- inner products of position eigenvectors with momentum eigenvectors in different order are complex conjugates

$$\langle x|p\rangle = A e^{\frac{ipx}{\hbar}} \quad \langle x|p\rangle = (\langle p|x\rangle)^*$$

$$\langle p|x\rangle = A e^{-\frac{ipx}{\hbar}}$$

- wave function in terms of momentum - $\tilde{\Psi}(p)$ ~

wave function in terms of position - $\psi(x)$

never both - $\psi(x,p)$ X !!

- wavelength of particle with function $e^{\frac{ipx}{\hbar}}$ is

$$\lambda = \frac{2\pi\hbar}{p} \quad \text{as } p \uparrow \lambda \downarrow$$

wavelength since if we add it to position x it doesn't change the wavefunction let $x' = x + \lambda = x + \frac{2\pi\hbar}{p}$

$$e^{\frac{ipx'}{\hbar}} = e^{\frac{ipx + 2\pi\hbar}{\hbar}} = e^{\frac{ipx}{\hbar}} e^{2\pi i} = e^{\frac{ipx}{\hbar}} \quad \text{so w/out change}$$

- called a wavefunction since $e^{\frac{ipx}{\hbar}} \Rightarrow e^{ix}$ represents cos/sin waves

momentum p $P(p) = |\langle p | \Psi \rangle|^2$ $\tilde{\Psi}(p) = \langle p | \Psi \rangle$

- resolving identity - identity operator I can be written as

↳ discrete wavefunctions $I = \sum_i |i\rangle \langle i|$

↳ continuous wavefunctions $I = \int dx |x\rangle \langle x|$ or $\int dp |p\rangle \langle p|$

so using $\tilde{\Psi}(p) = \langle p | \Psi \rangle$ and $I = \int dx |x\rangle \langle x|$

we get $\tilde{\Psi}(p) I = \int dx \underbrace{\langle p | x \rangle}_{\text{multiplied out } \langle p | \Psi \rangle \text{ since } \int dx} \langle x | \Psi \rangle$ so $\tilde{\Psi}(p) = \int dx \langle p | x \rangle \langle x | \Psi \rangle$

$x P \Psi(x) = -i\hbar x \frac{d\Psi(x)}{dx}$ $P x \Psi(x) = -i\hbar \frac{d(x\Psi(x))}{dx}$ } product rule

commutator $[x, P] \Psi(x)$
= $-i\hbar x \frac{d\Psi(x)}{dx} - i\hbar \Psi(x)$

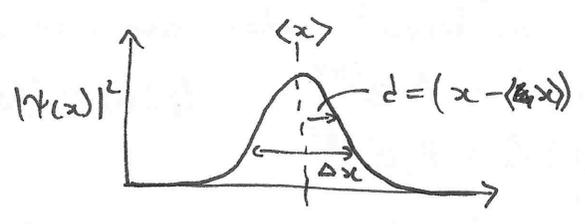
$[x, P] \Psi(x) = i\hbar x P \Psi(x) - P x \Psi(x)$
= $i\hbar \Psi(x)$

so $[x, P] = i\hbar$
since if it acts on $\Psi(x)$ it multiplies it by factor $i\hbar$

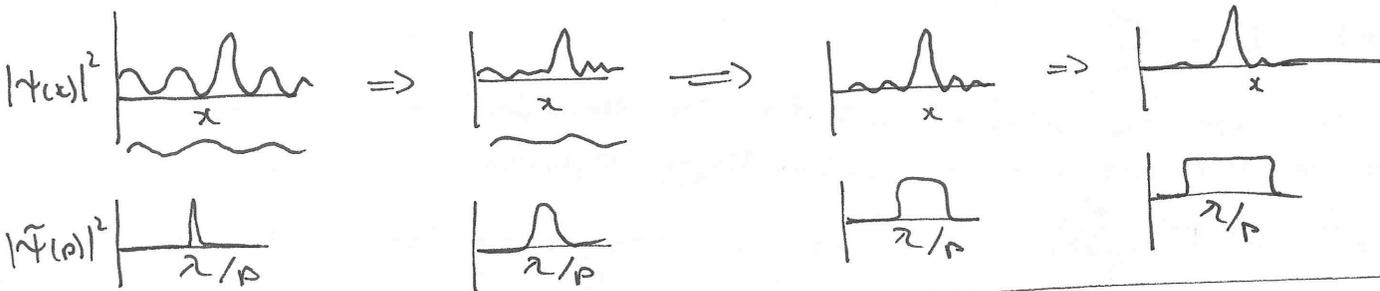
- Heisenberg's uncertainty principle
↳ since x, P don't commute $[x, P] \neq 0$

$$\begin{aligned} \Delta x \Delta p &\geq \frac{1}{2} |\langle \Psi | [x, P] | \Psi \rangle| \\ &\geq \frac{1}{2} |\langle \Psi | i\hbar | \Psi \rangle| \\ &\geq \frac{1}{2} |i\hbar \langle \Psi | \Psi \rangle| \\ &\geq \frac{\hbar}{2} \quad \text{since } \langle \Psi | \Psi \rangle = 1 \end{aligned}$$

Δx is measure of how spread out function is in relation to $\langle x \rangle$



- Heisenberg's uncertainty principle can be explored with Fourier transforms
 we superimpose waves of slightly different λ to get a single peak therefore know the $\langle x \rangle$ accurately but now we don't know the wavelength



- simple moving particle (non-relativistic)

on this eg simple Hamiltonian operator

$$H = cP = c \frac{d}{dx} - c i \hbar \frac{d}{dx}$$

plug into the independent Schrödinger equation

$$i \hbar \frac{\partial |\Psi\rangle}{\partial t} = H |\Psi\rangle$$

$$i \hbar \frac{\partial |\Psi\rangle}{\partial t} = -c i \hbar \frac{\partial |\Psi\rangle}{\partial x} \quad \text{or} \quad i \hbar \frac{\partial \psi(x,t)}{\partial t} = -c i \hbar \frac{\partial \psi(x,t)}{\partial x}$$

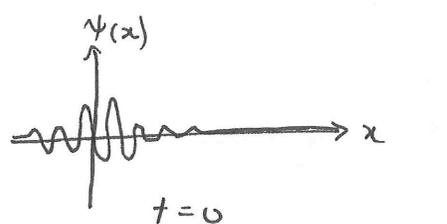
which gives: $\frac{\partial \psi(x,t)}{\partial t} = -c \frac{\partial \psi(x,t)}{\partial x}$

solution is $\psi(x,t) = x - ct$

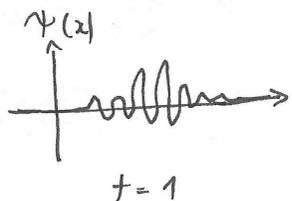
we have to normalize it

$$\int_{-\infty}^{\infty} \psi^*(x) \psi(x) dx$$

only in terms of space



at t+1



so wavepacket moved to the right

- classical Hamiltonian $H = \frac{1}{2}mv^2 = \frac{p^2}{2m}$ (free \therefore no external energy) so potential E term = 0

to quantize it replace $p \Rightarrow \hat{P}$

$$H = \frac{P^2}{2m} \quad P = -i \hbar \frac{\partial}{\partial x}$$

$$= (-i \hbar \frac{\partial}{\partial x}) (-i \hbar \frac{\partial}{\partial x}) \frac{1}{2m}$$

$$= \frac{\hbar^2}{2m} \frac{\partial^2}{\partial x^2}$$

substituting into $i \hbar \frac{\partial \psi}{\partial t} = H \psi$ gives:

$$i \hbar \frac{\partial \psi}{\partial t} = \frac{\hbar^2}{2m} \frac{\partial^2 \psi}{\partial x^2}$$

$$i \frac{\partial \psi}{\partial t} = \frac{\hbar}{2m} \frac{\partial^2 \psi}{\partial x^2}$$

this wave function spreads out over time



but λ is constant as long as a velocity so P is constant

- schrodinger's time independent ~~wave function~~ equation - $H|\psi\rangle = E|\psi\rangle$

so E is eigenvalue of hamiltonian

substituting we get $-\frac{\hbar^2}{2m} \frac{\partial^2 \psi(x)}{\partial x^2} = E\psi(x)$ assume solution $\psi(x) = e^{\frac{iPx}{\hbar}}$

we get $E = \frac{p^2}{2m}$

- now we can use this solution with the time dependant one to get the solution to the time dependant schrodinger equation

$\psi(x,t) = e^{\frac{i(px - \frac{p^2t}{2m})}{\hbar}}$ and any solution is sum or integral of these solutions in terms of p

$\psi(x,t) = \int \tilde{\psi}(p) \left(e^{\frac{i(px - \frac{p^2t}{2m})}{\hbar}} \right) dp$ also for wavefunction in terms of momentum $\tilde{\psi}(p,t) = \tilde{\psi}(p) e^{\frac{i(px - \frac{p^2t}{2m})}{\hbar}}$

- velocity and momentum

- velocity = time derivative of average position $v = \frac{d}{dt} \langle \psi | X | \psi \rangle$
in terms of wave functions $v = \frac{d}{dt} \int \psi^*(x,t) x \psi(x,t) dx$

- we know time derivative of expectation value of observable L is equal to $\frac{i}{\hbar}$ times expectation value of commutator of the hamiltonian H with the observable L $\frac{d}{dt} \langle L \rangle = \frac{i}{\hbar} \langle [H, L] \rangle$

- using hamiltonian from before $v = \frac{i}{2m\hbar} \langle [P^2, X] \rangle$

$[P^2, X] = P[P, X] + [P, X]P$ using $[P, X] = -i\hbar$ which we know
 $= -2i\hbar \langle P \rangle$

$v = \frac{2\hbar \langle P \rangle}{2m\hbar} = \frac{\langle P \rangle}{m}$ so $\langle P \rangle = mv$

- Quantization

1.) start with classical system - coordinates x_i and momenta p_i
- hamiltonian equation $H = \dots$

2.) replace classical phase space with linear vector space
state of states is represented in terms of coordinates by $\psi(x)$
or momenta by $\psi(p)$

3.) replace x_i and p_i in hamiltonian by X_i and P_i which act $X_i \psi = x_i \psi$

4.) hamiltonian is now an operator and can be used $P_i = -\hbar \frac{\partial}{\partial x_i}$
in the time dependant/independant equations

time dependant - shows development of wavefunction over time

time independant - enables us to calculate eigenvalues + eigenvectors of the hamiltonian corresponding to the Energy

- Forces - force gives rise to potential term in hamiltonian
 force can be quantized as potential and becomes operator V
 which is added to the hamiltonian to be used $H = T + V$

$V|\psi\rangle = V(x)\psi(x)$ so operator V multiplies wave function by a function V

new hamiltonian is $H = \frac{p^2}{2m} + V(x)$ schrodgers: $i\hbar \frac{\partial \psi}{\partial t} = -\frac{\hbar^2}{2m} \frac{\partial^2 \psi}{\partial x^2} + V(x)\psi$

so affects the way ψ changes with time and E $E\psi = -\frac{\hbar^2}{2m} \frac{\partial^2 \psi}{\partial x^2} + V(x)\psi$

- multiplying by x and by $V(x)$ are both commutative so x and $V(x)$ commute
 $[x, V(x)] = 0$ so $V(x)$ doesn't affect the before shown commutative relationship and $\langle p \rangle = mv$ still applies

- classical newtons equations of motion $\frac{dp}{dt} = F$

to try and quantize it calculate $\frac{d}{dt} \langle p \rangle$ using commutation with H

$\frac{d}{dt} \langle p \rangle = \frac{i}{\hbar} \langle [H, P] \rangle = \frac{i}{2m\hbar} \langle [P^2, P] \rangle + \frac{i}{\hbar} \langle [V, P] \rangle$ commutator commutes with any function of itself so $[P^2, P] = 0$

$= \frac{i}{\hbar} \langle [V, P] \rangle$
 $= \frac{i}{\hbar} \left\langle \left[i\hbar \frac{dV(x)}{dx} \right] \right\rangle$ using relationship

$= - \left\langle \frac{dV(x)}{dx} \right\rangle$ $[V(x), P] = i\hbar \frac{dV(x)}{dx}$

so we get $\frac{d}{dt} \langle p \rangle = - \left\langle \frac{dV(x)}{dx} \right\rangle$

- note: ~~derivative with respect~~ function of average of x ($\langle x \rangle$) is not the same as average of function of x $f(\langle x \rangle) \neq \langle f(x) \rangle$!

- classical limit

- assume incoming wave packet with features of size δx smaller than uncertainty in position so $\delta x < \Delta x$

- if sharp peaks of $V(x)$ due to force exist on scale smaller than incoming wave packet it is scattered so wave function breaks up

uncertainty principle
 $\Delta x \Delta p \geq \frac{\hbar}{2}$
 $\Delta x m \Delta v \geq \frac{\hbar}{2}$
 $\Delta x \Delta v \geq \frac{\hbar}{2m}$
 or $\Delta x \geq \frac{\hbar}{2m \Delta v}$

- gaussian wave packets - $\Delta x \Delta p$ is as small as possible so $\Delta x \Delta p = \hbar$

- Path integrals:

action $A = \int_{t_1}^{t_2} L(x, \dot{x}) dt$ $L = T - V$ in classical mechanics
here we deal with trajectory between two points

- in quantum mechanics there is no sense in a path between two points
so we ask: given particle starts at (x_1, t_1) what is the amplitude (probability) that it will show up at (x_2, t_2)

lets call the amplitude $C(x_1, t_1, x_2, t_2)$ so $C_{1,2}$

initial state $|\psi(t_1)\rangle = |x_1\rangle$ which evolves $|\psi(t_2)\rangle = e^{-iH(t_2-t_1)}|x_1\rangle$

$$C_{1,2} = \langle x_2 | e^{-iH(t_2-t_1)} | x_1 \rangle \text{ and define } t = t_2 - t_1$$
$$= \langle x_2 | e^{-iHt} | x_1 \rangle$$

using identity operator $I = \int dx |x\rangle \langle x|$

$$= \int dx \langle x_2 | e^{-\frac{iHt}{2}} | x \rangle \langle x | e^{-\frac{iHt}{2}} | x_1 \rangle$$

power is $\frac{1}{2}$ since we square rooted it to divide it in to interval

- so for each tiny time interval ϵ

- we include a factor $e^{-i\epsilon H}$

and in interval it gives term $\langle x_i | e^{-i\epsilon H} | x_{i+1} \rangle$

- if we let $U(\epsilon) = e^{-i\epsilon H}$

so we write it as $C_{1,2} = \langle x_2 | U^N | x_1 \rangle$

$$= \langle x_2 | U U U U \dots | x_1 \rangle$$

~~the term N is called the amplitude of the given path so we write~~

$$= \int dx \langle x_2 | e^{-i\epsilon H} | x \rangle \langle x | e^{-i\epsilon H} | x \rangle \langle x | e^{-i\epsilon H} | x_1 \rangle \dots$$

$$e^{\frac{iA}{\hbar}}$$

- A is the action for individual paths $e^{\frac{iA}{\hbar}}$ $C_{1,2} = \int_{\text{paths}} e^{\frac{iA}{\hbar}}$

so amplitude is integral over all possible paths between end points at infinitesimal time intervals $\epsilon \rightarrow 0$

↳ important for QED so remember.

- classical equation for harmonic oscillator

$$H = \frac{1}{2} \dot{x}^2 + \frac{1}{2} \omega^2 x^2 \quad \text{canonical momentum} = p = \frac{\partial L}{\partial \dot{x}} = \dot{x}$$

$$\text{so we get } H = \frac{1}{2} p^2 + \frac{1}{2} \omega^2 x^2$$

- quantum harmonic oscillator:

classical $H = \frac{1}{2} p^2 + \frac{1}{2} \omega^2 x^2$ $X|\psi(x)\rangle = x\psi(x)$

$H = \frac{1}{2} \left(-i\hbar \frac{\partial}{\partial x} \left(-i\hbar \frac{\partial}{\partial x} \psi(x) \right) \right) + \frac{1}{2} \omega^2 x^2 \psi(x)$ $P|\psi(x)\rangle = -i\hbar \frac{d}{dx} \psi(x)$

$H\psi\rangle = -\frac{\hbar^2}{2} \frac{\partial^2 \psi(x)}{\partial x^2} + \frac{1}{2} \omega^2 x^2 \psi(x)$

- substituting into schrodinger equation $i \frac{\partial \psi}{\partial t} = \frac{1}{\hbar} H\psi$

we get $i \frac{\partial \psi}{\partial t} = -\frac{\hbar}{2} \frac{\partial^2 \psi}{\partial x^2} + \frac{1}{2\hbar} \omega^2 x^2 \psi$

- substituting into $H|\psi_E\rangle = E|\psi_E\rangle$ we can find E levels

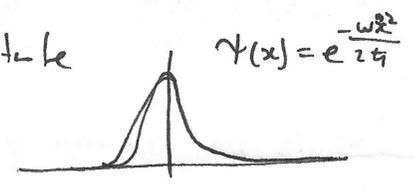
$-\frac{\hbar^2}{2} \frac{\partial^2 \psi_E(x)}{\partial x^2} + \frac{1}{2} \omega^2 x^2 \psi_E(x) = E \psi_E(x)$

- most values of E (solutions) grow exponentially as $x \rightarrow \infty$ or $-\infty$ so make no physical sense \therefore we create condition:

↳ physical solutions to schrodinger equation must be normalizable so filters out majority of E except few special ones

- lowest E state for harmonic oscillator = ground state

- ground state: - has no nodes
- has no zeroes for any potential



ground state wave function $\psi_0(x) = e^{-\frac{\omega}{2\hbar} x^2}$

substituting into schrodinger equation: $\frac{\hbar}{2} \omega e^{-\frac{\omega}{2\hbar} x^2} = E e^{-\frac{\omega}{2\hbar} x^2}$

ground state function $\psi_0(x)$ is the gaussian. so $E_0 = \frac{\omega\hbar}{2}$

- constructing annihilation/creation operators

hamiltonian in terms of operators $H = \frac{P^2 + \omega^2 X^2}{2}$ using: $a^2 + b^2 = (a+ib)(a-ib)$

but since X, P don't commute we can't just expand it so multiplying out we find we are off by factor of $\frac{\omega\hbar}{2}$

$= \frac{1}{2} (P+i\omega X)(P-i\omega X)$
actually $= \frac{1}{2} (P+i\omega X)(P-i\omega X) + \frac{\omega\hbar}{2}$

now define creation operator as $a^+ = \frac{-i}{\sqrt{2\omega\hbar}} (P+i\omega X)$

annihilation operator as $a^- = \frac{i}{\sqrt{2\omega\hbar}} (P-i\omega X)$

so hamiltonian becomes: $H = \omega\hbar (a^+ a^- + \frac{1}{2})$

and their commutator property $[a^+, a^-] = 1$

$[a^-, a^+] = -1$

- now define operator $N = a^+ a^-$ - number operator

$$H = \omega \hbar (N + \frac{1}{2})$$

for N we get commutator relationships: $[a^-, a^+] = 1$
 $[a^-, N] = a^-$
 $[a^+, N] = -a^+$

- create eigenvalue/value

$$N |u\rangle = u |u\rangle$$

$$\text{and act on it with } a^+ \quad N(a^+ |u\rangle) = [a^+ N - (a^+ N - N a^+)] |u\rangle$$

$$\left. \begin{array}{l} \text{since } |u\rangle \text{ is eigenvalue of } N \text{ with value } u \text{ we can} \\ \text{replace } (N+1) \text{ with } (u+1) \\ \text{if acting on } |u\rangle \end{array} \right\} \begin{array}{l} = a^+ (N+1) |u\rangle \\ = (u+1) (a^+ |u\rangle) \\ \underline{a^+ |u\rangle = |u+1\rangle} \end{array}$$

so we found a new eigen vector with eigenvalue increased by 1
so there is ∞ number of eigen vectors above it: $|u+2\rangle, |u+3\rangle \dots$

annihilation operator does opposite $a^- |u\rangle = |u-1\rangle$

$$\underline{a^+ |u\rangle = |u+1\rangle} \quad \underline{a^- |u\rangle = |u-1\rangle}$$

but we can't get below ground state to -ve E so there must be an eigen vector on which if a^- acts it gives back the same eigen vector so all lower ones are the same $a^- |0\rangle = 0$

this is the ground state $|0\rangle$ is not the 0 vector

- this quantizes E of harmonic oscillator to $E_n = \omega \hbar (n + \frac{1}{2})$
 $= \omega \hbar (\frac{1}{2}, \frac{3}{2}, \frac{5}{2}, \frac{7}{2} \dots)$

- $a^- |0\rangle = 0$ in terms of ground state $\psi_0(x)$:

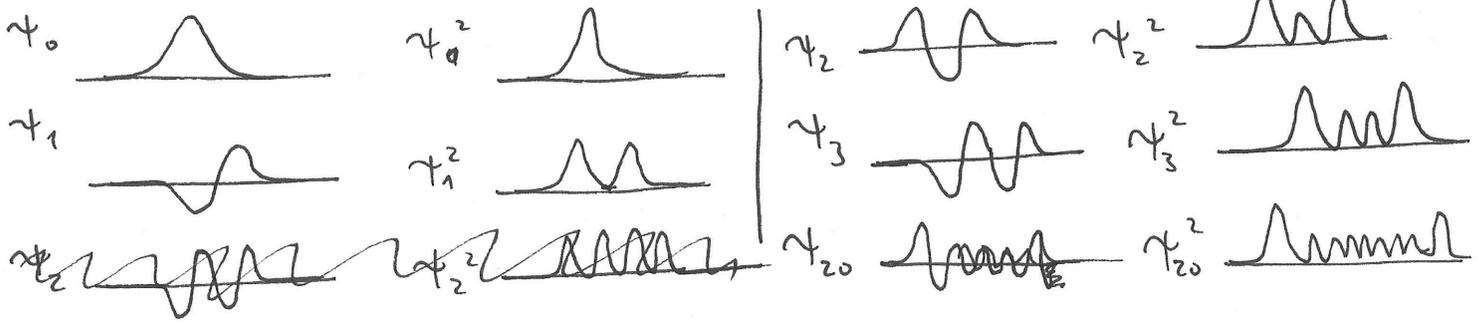
$$\frac{i}{\sqrt{2\omega\hbar}} (P - i\omega X) \psi_0(x) = 0 \quad \left\{ \begin{array}{l} \text{since } \frac{i}{\sqrt{2\omega\hbar}} \text{ is const we can divide by it} \\ (P - i\omega X) \psi_0(x) = 0 \end{array} \right.$$

$$\frac{d\psi_0}{dx} = -\frac{\omega x}{\hbar} \psi_0 \quad \left\{ \begin{array}{l} \text{replacing } P \text{ with } -i\hbar \frac{d}{dx} \\ \text{first order differential eqn} \end{array} \right.$$

solution is the gaussian $\psi_0 = e^{-\frac{\omega}{2\hbar} x^2}$

$$\begin{aligned} - \psi_1(x) &= (P + i\omega X) \psi_0(x) \\ &= (-i\hbar \frac{\partial}{\partial x} + i\omega x) e^{-\frac{\omega}{2\hbar} x^2} \\ &= i(\omega x - \hbar \frac{\partial}{\partial x}) e^{-\frac{\omega}{2\hbar} x^2} = i(\omega x e^{-\frac{\omega}{2\hbar} x^2} - \hbar \frac{\partial}{\partial x} (e^{-\frac{\omega}{2\hbar} x^2})) \\ &= i(\omega x e^{-\frac{\omega}{2\hbar} x^2} + \hbar 2x \frac{\omega}{2\hbar} e^{-\frac{\omega}{2\hbar} x^2}) \\ &= 2i\omega x e^{-\frac{\omega}{2\hbar} x^2} \\ \text{if we go on we get polynomials in terms of } x \text{ due to rule of } \frac{d}{dx} e^{ux} &\rightarrow \psi_1 = 2i\omega x \psi_0 \end{aligned}$$

- harmonic oscillator eigen functions



- $\omega = \frac{2\pi c}{\lambda}$ but $E \propto f$ so E in length of wave λ has to be $(n + \frac{1}{2}) h\omega$ $\frac{1}{2} h\omega$ - zero point E

- ignoring zero point E so $E_n = \frac{2\pi h c}{\lambda} n$ $n \in +\mathbb{Z}$ or 0

so E is quantized in electromagnetic wave

$$E(\lambda) = \frac{2\pi h c}{\lambda}$$

↳ energy of a single photon